

## Standard 3 gamma spectrometer corrections (T52)

Standard 3 Corrections performs the following corrections to the standard gamma-ray spectrometer data channels Potassium, Uranium, Thorium and Total Count

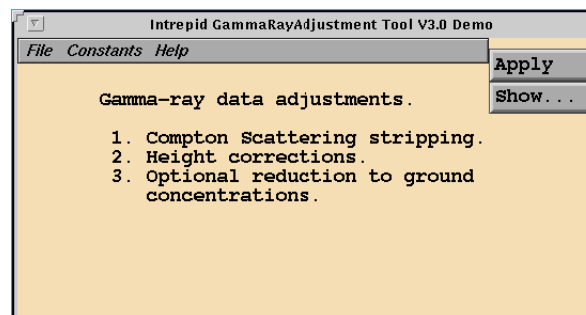
- 1 Normalising measurements to a standard distance above ground level
- 2 Stripping correction for Compton scattering
- 3 Optional) Conversion of count rates to elemental ground concentrations.

See "[Standard 3 corrections theory](#)" in [Radiometrics theoretical background \(R25\)](#) for this tool's background theory.

### Using the S3GSC corrections tool

>> *To use S3GSC with the INTREPID graphic user interface*

- 1 Choose Standard 3 from the Radiometrics menu in the Project Manager, or use the command gamadj.exe. INTREPID displays the S3GSC window



- 2 If you have previously prepared file specifications and parameter settings for S3GSC, load the corresponding task specification file using Load Options from the File menu. (See [Specifying input and output files](#) for detailed instructions.) If all of the specifications are correct in this file, go to step 7. If you wish to modify any settings, carry out the following steps as required.
- 3 Specify the dataset to be processed. Use Specify Radiometrics Dataset from the File menu. (See [Specifying input and output files](#) for detailed instructions.)
- 4 Specify the atmospheric temperature and pressure files. Use Open Air Input Database from the File menu. (See [Specifying input and output files](#) for detailed instructions.)
- 5 Specify the line type field if required for eliminating selected traverse lines from the process. Use LineType from the File menu. (See [Specifying input and output files](#) for detailed instructions.)
- 6 Specify the parameters for the corrections using the options from the Constants menu (See [Standard 3 correction constants](#) for details).
- 7 When you have made specifications and settings according to your requirements, choose Apply. INTREPID will perform the corrections. The corrected data overwrites the original data.
- 8 If you wish to record the specifications for this process in a **.job** file in order to repeat a similar task later or for some other reason, use Save Options from the file menu. (See [Specifying input and output files](#) for detailed instructions.)
- 9 If you wish to repeat the process, repeat steps 2–7, varying the specifications as required.

**10** To exit from S3GSC, choose Quit from the File menu.

To view your current file specifications and parameter settings choose Show. INTREPID displays them in a separate window. See [Displaying options and using Task specification files](#) for details and an example of a set of specifications

S3GSC performs stripping before height correction. The coefficients would be different if it performed the corrections in the reverse order.

In the current version there is no graphic display of the S3GSC process.

You can view Help information by choosing options from the Help menu (See [Help](#)).

You can execute S3GSC as a batch task by command using a task specification (`.job`) file that you have previously prepared. See [Displaying options and using Task specification files](#) for details.

## Specifying input and output files

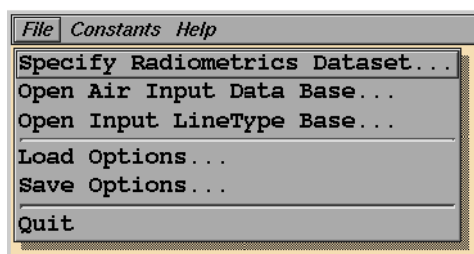
To use S3GSC you must specify

- The line dataset to be processed, with aliases to identify the required fields,
- The dataset containing the air temperature and pressure readings,
- The traverse line type field.

To identify the fields in the dataset, give them aliases and specify the aliases and the file names in the dataset. See "[Vector dataset field aliases](#)" in [INTREPID database, file and data structures \(R05\)](#) for more information about aliases.

This section lists the aliases required and has instructions for specifying the files for the S3GSC process.

Choose the options as required from the File menu.



In each case INTREPID displays an Open dialog box. Use the directory and file selector to locate the line dataset directory or file you require. (See "[Specifying input and output files](#)" in [Introduction to INTREPID \(R02\)](#) for information about specifying files).

**Specify Radiometrics Dataset** Use this option to specify the lines dataset you wish to process.

To perform the S3GSC corrections, you must specify the following field aliases:

Alias	Field
Potassium	potassium channel data (counts/second)
Uranium	uranium channel data (counts/second)
Thorium	thorium channel data (counts/second)
TotalCount	total count data (counts/second)
Clearance	height of the aircraft above ground level (metres)
Fiducial	fiducial record
SurveyNumber	(optional) survey number field for linking to the temperature/pressure data
FlightNumber	(optional) flight number field for linking to the temperature/pressure data
LineNumber	traverse line number field for linking to the temperature/pressure data

When you choose Apply, INTREPID will modify the fields with aliases **Potassium**, **Uranium**, **Thorium** and **TotalCount** in the line dataset you specify.

**Open Air Input Dataset** Use this to specify the line dataset that contains the atmospheric pressure and temperature data. This data may have a lower sampling rate than the radiometrics data, and therefore may be in a different INTREPID dataset.

INTREPID will only adjust the height data for temperature and pressure if you specify this dataset. For INTREPID to use this data you must specify the following aliases for fields

Alias	Field
Temperature	temperature (degrees Celsius)
Pressure	atmospheric pressure (millibars)
Fiducial	fiducial data
SurveyNumber	(optional) the survey number field for linking to the radiometrics data
FlightNumber	(optional) the flight number field for linking to the radiometrics data
LineNumber	traverse line number field for linking to the radiometrics data

Note that if the temperature and pressure data is in a different line dataset, the aliases must be listed in the actual line dataset that contains the temperature and pressure data.

**Open LineType** Use this to specify the line type field for eliminating selected traverse lines from the process. If you do not specify a line type field INTREPID will process all traverse lines data. If you specify a line type field, INTREPID will process all traverse lines except those whose line type is 16384.

**Load Options** If you wish to use an existing task specification file to specify the S3GSC process, use this menu option to specify the task specification file required. INTREPID will load the file and use its contents to set all of the parameters for the S3GSC process. (See [Displaying options and using Task specification files](#) for information about task specification files).

**Save Options** If you wish to save the current S3GSC file specifications and parameter settings as an task specification file, use this menu option to specify the filename and save the file. (See [Displaying options and using Task specification files](#) for information about task specification files).

### Sample aliases

#### Alias Begin

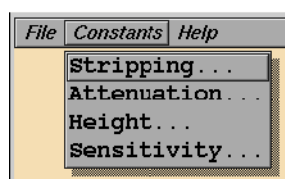
```
Potassium = potassium
Uranium = uranium
Thorium = thorium
TotalCount = totalcount
Fiducial = FID
Clearance = clearance
LineNumber = LINE
FlightNumber = flight
SurveyNumber = survey
```

#### Alias End

## Standard 3 correction constants

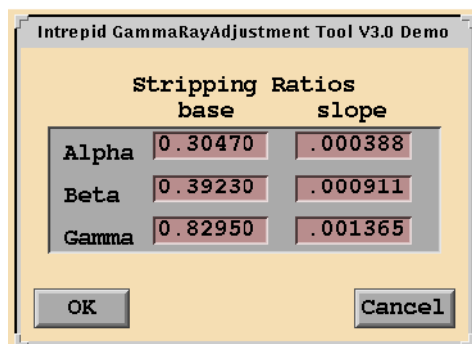
Use the options from the Constants menu to define

- The stripping, height, attenuation, and sensitivity constants,
- The height above which data is to be ignored due to the unreliability of the height attenuation correction,
- The nominal survey height to which you require the data reduced.



## The Stripping Ratios

To set the Stripping Ratios, choose Stripping from the Constants menu. INTREPID displays the Stripping Ratios dialog box.



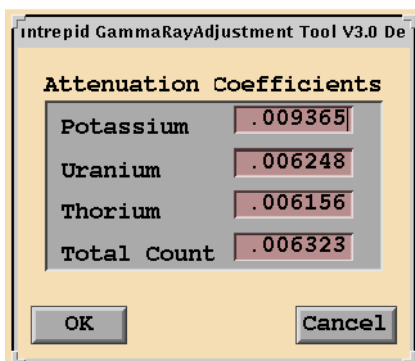
	base	slope
Alpha	0.30470	.000388
Beta	0.39230	.000911
Gamma	0.82950	.001365

Use the dialog box to set the base level (ground level value) and slope (change per metre) of the stripping ratios  $\alpha$ ,  $\beta$ ,  $\gamma$  in equations (1) above. (See ["Compton scattering correction \(stripping\)—theory" in Radiometrics theoretical background \(R25\)](#)).

INTREPID will set default values according to the AGSO system (See the footnote to ["Compton scattering correction \(stripping\)—theory" in Radiometrics theoretical background \(R25\)](#)).

## The Attenuation Coefficients

To set the Attenuation Coefficients, choose Attenuation from the Constants menu. INTREPID displays the Attenuation Coefficients dialog box.



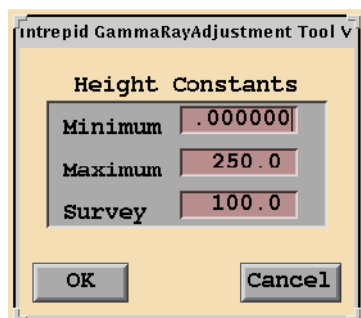
Potassium	.009365
Uranium	.006248
Thorium	.006156
Total Count	.006323

Use the dialog box to set the Attenuation Coefficients (values of  $u$  for Potassium Uranium Thorium and Total Count in equation (2) above (See ["Height correction theory" in Radiometrics theoretical background \(R25\)](#)).

INTREPID will set default values according to the AGSO system (See the footnote to ["Height correction theory" in Radiometrics theoretical background \(R25\)](#))

## Height Constants

To set the Height Constants, choose Height from the Constants menu. INTREPID displays the Height Constants dialog box.



Use the dialog box to set the Height Constants.

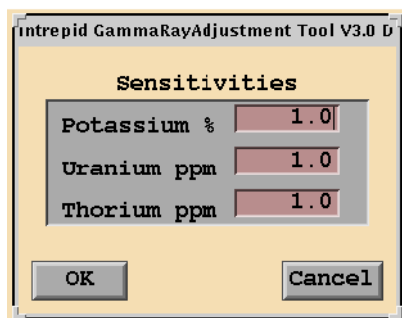
**minimum height** the height below which you wish INTREPID to ignore the radiometrics data (default value 0).

**maximum height** the height above which you wish INTREPID to ignore the radiometrics data due to the unreliability of the height attenuation correction. (default value 250m).

**survey** nominal height to which you wish to reduce your survey data (H in equation (2) (See "[Height correction theory](#)" in [Radiometrics theoretical background \(R25\)](#)) (default value 100m).

## Sensitivities

To set the Sensitivities, choose Sensitivity from the Constants menu. INTREPID displays the Sensitivities dialog box.



Use this dialog box to specify the factors for converting count rates to ground concentrations for Potassium, Uranium, and Thorium. Express each factor in counts per second per unit concentration. For each element INTREPID will divide the measured and corrected count rates by the factor you specify to obtain the ground concentrations of

- Potassium (in percentage by mass),
- Uranium (in parts per million) and
- Thorium (in parts per million)

If you do not have values for sensitivity constants or do not wish to convert your data to ground concentration, set the values of the sensitivity constants to 1.0. INTREPID will leave your data in counts/second.

See "[Sensitivity conversions—theory](#)" in [Radiometrics theoretical background \(R25\)](#) for more information about sensitivity.

If you do not set a value for the sensitivity for one or more elements, INTREPID will use the default value **1.0** for each unspecified constant.

## Apply

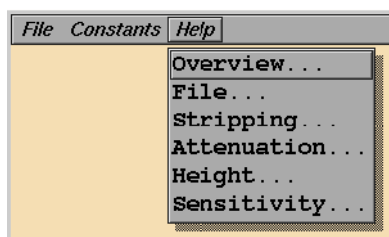
When you choose Apply, INTREPID performs the corrections and modifies the files indicated by the **Potassium**, **Uranium** and **Thorium** aliases.

## Exit

To exit from S3GSC choose Quit from the file menu.

## Help

You can use the help menu to display help text on the topics shown in the menu illustration below.



## Displaying options and using Task specification files

### Displaying options

To display the current file specifications and parameter settings

Choose Show in the S3GSC window. INTREPID displays the current options in a separate report window.

### Using task specification files

You can store sets of file specifications and parameter settings for S3GSC in task specification (**.job**) files.

#### >> *To create a task specification file with the S3GSC tool*

- 1 Specify all files and parameters.
- 2 If possible, execute the task (choose Apply) to ensure that it will work.
- 3 Choose Save Options from the File menu. Specify a task specification file (INTREPID will add the extension **.job**) INTREPID will create the file with the settings current at the time of the Save Options operation.

For full instructions on creating and editing task specification files see Appendix F Task Specification Files in Volume 1

#### >> *To use a task specification file in an interactive S3GSC session*

Load the task specification (**.job**) file (File menu, Load Options), modify any settings as required, then choose Apply.

#### >> *To use a task specification file for a batch mode S3GSC task*

Type the command **gamadj.exe** with the switch **-batch** followed by the name of the task specification file.

For example, if you had a task specification file called **surv329.job** you would use the command

```
gamadj.exe -batch surv329.job
```



## Task specification file example

Here is an example of a S3GSC task specification file.

```
Process Begin
  Name = Input??
  Parameters Begin
    RadiometricsDataBaseDirectory = "/disk1/survey/rad/SEC"
    AirDataBaseDirectory = "/disk1/survey/rad/TEN"
    alphaBase = 0.46
    alphaSlope = 0.00
    betaBase = 0.44
    betaSlope = 0.00
    gammaBase = 0.81
    gammaSlope = 0.00
    PotassiumAttenuation = 0.01
    UraniumAttenuation = 0.01
    ThoriumAttenuation = 0.01
    TotalCountAttenuation = 0.01
    MinimumHeight = 0.00
    MaximumHeight = 250.00
    NominalHeight = 150.00
    PotassiumSensitivity = 1.00
    UraniumSensitivity = 1.00
    ThoriumSensitivity = 1.00
    UseTempAndPressure = Yes
  Parameters End
Process End
```